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ПП

10. Sequential Data

Bayes Filter (Rep.)

We can describe the overall process using a Dynamic Bayes Network:



• This incorporates the following Markov assumptions:

$$p(z_t \mid x_{0:t}, u_{1:t}, z_{1:t}) = p(z_t \mid x_t) \text{ (measurement)}$$

$$p(x_t \mid x_{0:t-1}, u_{1:t}, z_{1:t}) = p(x_t \mid x_{t-1}, u_t) \text{ (state)}$$



Bayes Filter Without Actions

Removing the action variables we obtain:



• This incorporates the following Markov assumptions:

$$p(z_t \mid x_{0:t}, \quad z_{1:t}) = p(z_t \mid x_t) \text{ (measurement)}$$

$$p(x_t \mid x_{0:t-1}, \quad z_{1:t}) = p(x_t \mid x_{t-1}) \text{ (state)}$$
Machine Learning for



 Observations in sequential data should not be modeled as independent variables such as:



- Examples: weather forecast, speech, handwritten text, etc.
- The observation at time t depends on the observation(s) of (an) earlier time step(s):







• The joint distribution is therefore (d-sep):

$$p(\mathbf{z}_1 \dots \mathbf{z}_n) = p(\mathbf{z}_1) \prod_{i=2}^n p(\mathbf{z}_i \mid \mathbf{z}_{i-1})$$

 However: often data depends on several earlier observations (not just one)







- Problem: number of stored parameters grows exponentially with the order of the Markov chain
- Question: can we model dependency of all previous observations with a limited number of parameters?



Idea: Introduce hidden (unobserved) variables:







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Now we have: dsep($\mathbf{x}_n, \{\mathbf{x}_1, \dots, \mathbf{x}_{n-2}\}, \mathbf{x}_{n-1}$) $\Leftrightarrow p(\mathbf{x}_n \mid \mathbf{x}_1, \dots, \mathbf{x}_{n-2}, \mathbf{x}_{n-1}) = p(\mathbf{x}_n \mid \mathbf{x}_{n-1})$ But: $\neg dsep(\mathbf{z}_n, \{\mathbf{z}_1, \dots, \mathbf{z}_{n-2}\}, \mathbf{z}_{n-1})$ $\Leftrightarrow p(\mathbf{z}_n \mid \mathbf{z}_1, \dots, \mathbf{z}_{n-2}, \mathbf{z}_{n-1}) \neq p(\mathbf{z}_n \mid \mathbf{z}_{n-1})$

And: number of parameters is nK(K-1) + const.

Machine Learning for Computer Vision



Example

- Place recognition for mobile robots
- 3 different states: corridor, room, doorway
- Problem: misclassifications
- Idea: use information from previous time step





General Formulation of an HMM

- 1.Discrete random variables
 - **Observation** variables: $\{z_n\}, n = 1..N$
 - Discrete **state** variables (unobservable): $\{x_n\}, n = 1..N$
 - **Number** of states $K: x_n \in \{1...K\}$

2. Transition model $p(x_i | x_{i-1})$

- Markov assumption (x_i only depends on x_i)
- Represented as a *K*×*K* transition matrix *A*
- Initial probability: $p(x_0)$ repr. as π_1, π_2, π_3

3. Observation model $p(z_i|x_i)$ with parameters φ

- Observation only depends on the current state
- Example: output of a "local" place classifier



Model Parameters

θ

The Trellis Representation





Application Example (1)

- Given an observation sequence $z_1, z_2, z_3...$
- Assume that the model parameters $\theta = (A, \pi, \phi)$ are known
- What is the probability that the given observation sequence is actually observed under this model,
 i.e. the data likelihood p(Z| θ)?
- If we are given several different models, we can choose the one with highest probability
- Expressed as a supervised learning problem, this can be interpreted as the inference step (classification step)





Application Example (2)

Based on the data likelihood we can solve two different kinds of problems:

- Filtering: computes $p(\mathbf{x}_n | \mathbf{z}_{1:n})$, i.e. state probability only based on previous observations
- Smoothing: computes p(x_n | z_{1:N}), state probability based on all observations (including those from the future)







Application Example (3)

- Given an observation sequence $z_1, z_2, z_3...$
- Assume that the model parameters $\theta = (A, \pi, \varphi)$ are known
- What is the state sequence x₁, x₂, x₃... that
 explains best the given observation sequence?
- In the case of place recognition: which is the sequence of truly visited places that explains best the sequence of obtained place labels (classifications)?





Application Example (4)

- Given an observation sequence $z_1, z_2, z_3...$
- What are the optimal model parameters $\theta = (A, \pi, \phi)$?
- This can be interpreted as the training step
- It is in general the most difficult problem



Summary: 4 Operations on HMMs

- 1. Compute data likelihood $p(Z|\theta)$ from a known model
 - Can be computed with the forward algorithm
- 2. Filtering or Smoothing of the state probability
 - Filtering: forward algorithm
 - Smoothing: forward-backward algorithm
- 3. Compute optimal state sequence with a known model
 - Can be computed with the Viterbi-Algorithm
- 4. Learn model parameters for an observation sequence
 - Can be computed using Expectation-Maximization (or Baum-Welch)





Goal: compute $p(Z|\theta)$ (we drop θ in the following)

$$p(\mathbf{z}_1,\ldots,\mathbf{z}_n) = \sum_{\mathbf{x}_n} p(\mathbf{z}_1,\ldots,\mathbf{z}_n,\mathbf{x}_n) =: \sum_{\mathbf{x}_n} \alpha(\mathbf{x}_n)$$





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We can calculate α recursively:

$$\alpha(\mathbf{x}_n) = p(\mathbf{z}_n \mid \mathbf{x}_n) \sum_{\mathbf{x}_{n-1}} \alpha(\mathbf{x}_{n-1}) p(\mathbf{x}_n \mid \mathbf{x}_{n-1})$$



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This is the same recursive formula as we had in the first lecture!



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Filtering:
$$p(\mathbf{x}_n \mid \mathbf{z}_1, \dots, \mathbf{z}_n) = \frac{p(\mathbf{z}_1, \dots, \mathbf{z}_n, \mathbf{x}_n)}{p(\mathbf{z}_1, \dots, \mathbf{z}_n)} = \frac{\alpha(\mathbf{x}_n)}{\sum_{\mathbf{x}_n} \alpha(\mathbf{x}_n)}$$



The Forward-Backward Algorithm

- As before we set $\alpha(\mathbf{x}_n) = p(\mathbf{z}_1, \dots, \mathbf{z}_n, \mathbf{x}_n)$
- We also define $\beta(\mathbf{x}_n) = p(\mathbf{z}_{n+1}, \dots, \mathbf{z}_N \mid \mathbf{x}_n)$

e.g. *n* = 5:





The Forward-Backward Algorithm

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- We also define $\beta(\mathbf{x}_n) = p(\mathbf{z}_{n+1}, \dots, \mathbf{z}_N \mid \mathbf{x}_n)$
- This can be recursively computed (backwards):

 $\beta(\mathbf{x}_{n-1}) = p(\mathbf{z}_n, \dots, \mathbf{z}_N \mid \mathbf{x}_{n-1})$ $= \sum p(\mathbf{x}_n, \mathbf{z}_n, \dots, \mathbf{z}_N \mid \mathbf{x}_{n-1})$ $= \sum p(\mathbf{z}_{n+1},\ldots,\mathbf{z}_N \mid \mathbf{x}_n,\mathbf{y}_n,\mathbf{y}_{n-1})p(\mathbf{x}_n,\mathbf{z}_n \mid \mathbf{x}_{n-1})$ X_n $= \sum p(\mathbf{z}_{n+1},\ldots,\mathbf{z}_N \mid \mathbf{x}_n) p(\mathbf{z}_n \mid \mathbf{x}_{n-1},\mathbf{x}_n) p(\mathbf{x}_n \mid \mathbf{x}_{n-1})$ $= \sum \beta(\mathbf{x}_n) p(\mathbf{z}_n \mid \mathbf{x}_n) p(\mathbf{x}_n \mid \mathbf{x}_{n-1})$ \mathbf{X}_n



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- We also define $\beta(\mathbf{x}_n) = p(\mathbf{z}_{n+1}, \dots, \mathbf{z}_N \mid \mathbf{x}_n)$
- This can be recursively computed (backwards):

$$\beta(\mathbf{x}_n) = \sum_{\mathbf{x}_{n+1}} \beta(\mathbf{x}_{n+1}) p(\mathbf{z}_{n+1} \mid \mathbf{x}_{n+1}) p(\mathbf{x}_{n+1} \mid \mathbf{x}_n)$$

- This is exactly the same as the message-passing algorithm ("sum-product")!
 - forward messages α_n (vector of length *K*)
 - backward messages β_n (vector of length *K*)



Smoothing with Forward-Backward

First we compute $p(\mathbf{x}_n, \mathbf{z}_1, \dots, \mathbf{z}_N)$:

 $p(\mathbf{x}_n, \mathbf{z}_1, \ldots, \mathbf{z}_N) = p(\mathbf{z}_1, \ldots, \mathbf{z}_N \mid \mathbf{x}_n) p(\mathbf{x}_n)$

$$= p(\mathbf{z}_1, \dots, \mathbf{z}_n | \mathbf{x}_n) p(\mathbf{z}_{n+1}, \dots, \mathbf{z}_N | \mathbf{x}_n) p(\mathbf{x}_n)$$
$$= p(\mathbf{z}_1, \dots, \mathbf{z}_n, \mathbf{x}_n) p(\mathbf{z}_{n+1}, \dots, \mathbf{z}_N | \mathbf{x}_n)$$
$$= \alpha(\mathbf{x}_n) \beta(\mathbf{x}_n)$$



Smoothing with Forward-Backward

First we compute $p(\mathbf{x}_n, \mathbf{z}_1, \dots, \mathbf{z}_N)$:

 $p(\mathbf{x}_n, \mathbf{z}_1, \ldots, \mathbf{z}_N) = \alpha(\mathbf{x}_n)\beta(\mathbf{x}_n)$

with that we can compute $p(\mathbf{z}_1, \ldots, \mathbf{z}_N)$:

$$p(\mathbf{z}_1,\ldots,\mathbf{z}_N) = \sum_{\mathbf{x}_n} p(\mathbf{x}_n,\mathbf{z}_1,\ldots,\mathbf{z}_N) = \sum_{\mathbf{x}_n} \alpha(\mathbf{x}_n)\beta(\mathbf{x}_n)$$



Smoothing with Forward-Backward

First we compute $p(\mathbf{x}_n, \mathbf{z}_1, \dots, \mathbf{z}_N)$:

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with that we can compute $p(\mathbf{z}_1, \ldots, \mathbf{z}_N)$:

$$p(\mathbf{z}_1,\ldots,\mathbf{z}_N) = \sum_{\mathbf{x}_n} p(\mathbf{x}_n,\mathbf{z}_1,\ldots,\mathbf{z}_N) = \sum_{\mathbf{x}_n} \alpha(\mathbf{x}_n)\beta(\mathbf{x}_n)$$

and finally:

$$p(\mathbf{x}_n \mid \mathbf{z}_1, \dots, \mathbf{z}_N) = \frac{p(\mathbf{x}_n, \mathbf{z}_1, \dots, \mathbf{z}_N)}{p(\mathbf{z}_1, \dots, \mathbf{z}_N)} = \frac{\alpha(\mathbf{x}_n)\beta(\mathbf{x}_n)}{\sum_{\mathbf{x}_n} \alpha(\mathbf{x}_n)\beta(\mathbf{x}_n)}$$



2. Computing the Most Likely States

• Goal: find a state sequence $\mathbf{x}_1, \mathbf{x}_2, \mathbf{x}_3, \dots$ that maximizes the probability $p(X, Z | \theta)$

• Define
$$\delta(\mathbf{x}_n) = \max_{\mathbf{x}_1, \dots, \mathbf{x}_{n-1}} p(\mathbf{x}_1, \dots, \mathbf{x}_n \mid \mathbf{z}_1, \dots, \mathbf{z}_n)$$

This is the probability of state j by taking the most probable path.





2. Computing the Most Likely States

• Goal: find a state sequence $x_1, x_2, x_3...$ that maximizes the probability $p(X,Z|\theta)$

• Define
$$\delta(\mathbf{x}_n) = \max_{\mathbf{x}_1, \dots, \mathbf{x}_{n-1}} p(\mathbf{x}_1, \dots, \mathbf{x}_n \mid \mathbf{z}_1, \dots, \mathbf{z}_n)$$

This can be computed recursively: $\delta(\mathbf{x}_n) = \max_{\mathbf{x}_{n-1}} \delta(\mathbf{x}_{n-1}) p(\mathbf{x}_n | \mathbf{x}_{n-1}) p(\mathbf{z}_n, | \mathbf{x}_n)$

we also have to compute the argmax:

$$\psi(\mathbf{x}_n) = \arg \max_{\mathbf{x}_{n-1}} \delta(\mathbf{x}_{n-1}) p(\mathbf{x}_n \mid \mathbf{x}_{n-1}) p(\mathbf{z}_n, \mid \mathbf{x}_n)$$



The Viterbi algorithm

- Initialize:
 - $\delta(\mathbf{x}_0) = p(\mathbf{x}_0) p(\mathbf{z}_0 | \mathbf{x}_0)$
 - $\psi(\mathbf{x}_0) = 0$
- Compute recursively for *n*=1...*N*:
 - $\delta(\mathbf{x}_n) = p(\mathbf{z}_n | \mathbf{x}_n) \max_{\mathbf{x}_{n-1}} [\delta(\mathbf{x}_{n-1}) p(\mathbf{x}_n | \mathbf{x}_{n-1})]$
 - $\psi(\mathbf{x}_n) = \underset{x_{n-1}}{\operatorname{argmax}} \left[\delta(\mathbf{x}_{n-1}) p(\mathbf{x}_n | \mathbf{x}_{n-1}) \right]$
- On termination:
 - $p(Z, X|\theta) = \max_{x_N} \delta(x_N)$ • $x_N^* = \operatorname*{argmax}_{x_N} \delta(x_N)$
- Backtracking:

•
$$\mathbf{x}_n^* = \psi(\mathbf{x}_{n+1})$$



3. Learning the Model Parameters

- Given an observation sequence $z_1, z_2, z_3...$
- Find optimal model parameters $\theta = \pi, A, \varphi$
- We need to maximize the likelihood $p(Z|\theta)$
- Can not be solved in closed form
- Iterative algorithm: Expectation Maximization (EM) or for the case of HMMs: Baum-Welch algorithm



3. Learning the Model Parameters

Idea: instead of maximizing

$$p(\mathbf{z}_1,\ldots,\mathbf{z}_N \mid \theta) = \sum_X p(\mathbf{z}_1,\ldots,\mathbf{z}_N,\mathbf{x}_1,\ldots,\mathbf{x}_N \mid \theta)$$

• we maximize the expected log likelihood:

$$\sum_{X} p(\mathbf{x}_1, \dots, \mathbf{x}_N \mid \mathbf{z}_1, \dots, \mathbf{z}_N, \theta) \log p(\mathbf{z}_1, \dots, \mathbf{z}_N, \mathbf{x}_1, \dots, \mathbf{x}_N \mid \theta)$$

- it can be shown that this is a lower bound of the actual log-likelihood $p(Z|\theta)$
- this is exactly the Expectation-Maximization (EM) algorithm



- E-Step (assuming we know π , A, ϕ , i.e. θ^{old})
- Define the posterior probability of being in state i at step k:
- Define $\gamma(\mathbf{x}_n) = p(\mathbf{x}_n | Z)$





- E-Step (assuming we know π ,A, ϕ , i.e. θ^{old})
- Define the posterior probability of being in state i at step k:
- Define $\gamma(\mathbf{x}_n) = p(\mathbf{x}_n | \mathbf{z}_1, \dots, \mathbf{z}_n)$
- It follows that $\gamma(\mathbf{x}_n) = \alpha(\mathbf{x}_n) \beta(\mathbf{x}_n) / p(Z)$



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- It follows that $\gamma(\mathbf{x}_n) = \alpha(\mathbf{x}_n) \beta(\mathbf{x}_n) / p(Z)$
- Define $\xi(\mathbf{x}_{n-1},\mathbf{x}_n) = p(\mathbf{x}_{n-1},\mathbf{x}_n|Z)$
- It follows that

$$\xi(\mathbf{x}_{n-1},\mathbf{x}_n) = \alpha(\mathbf{x}_{n-1})p(\mathbf{z}_n|\mathbf{x}_n)p(\mathbf{x}_n|\mathbf{x}_{n-1})\beta(\mathbf{x}_n) / p(\mathbf{Z})$$



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- Define the posterior probability of being in state i at step k:
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- It follows that $\gamma(\mathbf{x}_n) = \alpha(\mathbf{x}_n) \beta(\mathbf{x}_n) / p(Z)$
- Define $\xi(x_{n-1}, x_n) = p(x_{n-1}, x_n | Z)$
- It follows that

 $\xi(\mathbf{x}_{n-1},\mathbf{x}_n) = \alpha(\mathbf{x}_{n-1})p(\mathbf{z}_n|\mathbf{x}_n)p(\mathbf{x}_n|\mathbf{x}_{n-1})\beta(\mathbf{x}_n) / p(\mathbf{Z})$

• The algorithm maximizes: $Q(\theta, \theta^{old}) = \sum p(X|Z, \theta^{old}) \log p(Z, X|\theta)$ "Expected complete data log-likelihood"



- Maximizing Q also maximizes the likelihood: $p(Z|\theta) \ge p(Z|\theta^{\text{old}})$
- M-Step:

$$\pi_k = \frac{\sum_{\mathbf{x}} \gamma(\mathbf{x}) x_{1k}}{\sum_{j=1} \sum_{\mathbf{x}} \gamma(\mathbf{x}) x_{1j}}$$

here, we need forward and backward step!

$$A_{jk} = \frac{\sum_{t=2}^{T} \xi(x_{t-1,j}, x_{tk})}{\sum_{l=1}^{K} \sum_{t=2}^{T} \xi(x_{t-1,j}, x_{tl})}$$

- With these new values, Q is recomputed
- This is done until the likelihood does not increase anymore (convergence)



The Baum-Welsh Algorithm - Summary

- Start with an initial estimate of $\theta = (\pi, A, \phi)$ e.g. uniformly and k-means for ϕ
- Compute $Q(\theta, \theta^{old})$ (E-Step)
- Maximize Q (M-step)
- Iterate E and M until convergence
- In each iteration one full application of the forward-backward algorithm is performed
- Result gives a local optimum
- For other local optima, the algorithm needs to be started again with new initialization



The Scaling problem

Probability of sequences



- Probabilities are very small
- The product of the terms soon is very small
- Usually: converting to log-space works
- But: we have sums of products!
- Solution: Rescale/Normalize the probability during the computation, e.g.:

$$\hat{\alpha}(\mathbf{x}_n) = \alpha(\mathbf{x}_n) / p(\mathbf{z}_1, \mathbf{z}_2, \dots, \mathbf{z}_n)$$



Summary

- HMMs are a way to model sequential data
- They assume discrete states
- Three possible operations can be performed with HMMs:
 - Data likelihood, given a model and an observation
 - Most likely state sequence, given a model and an observation
 - Optimal Model parameters, given an observation
- Appropriate scaling solves numerical problems
- HMMs are widely used, e.g. in speech recognition

